



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:01 AM UTC

PDB ID : 6N1G / pdb_00006n1g
Title : Crystal structure of Aquaglyceroporin AQP7
Authors : Vahedi-Faridi, A.; Lodowski, D.; Kowatz, T.
Deposited on : 2018-11-08
Resolution : 4.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

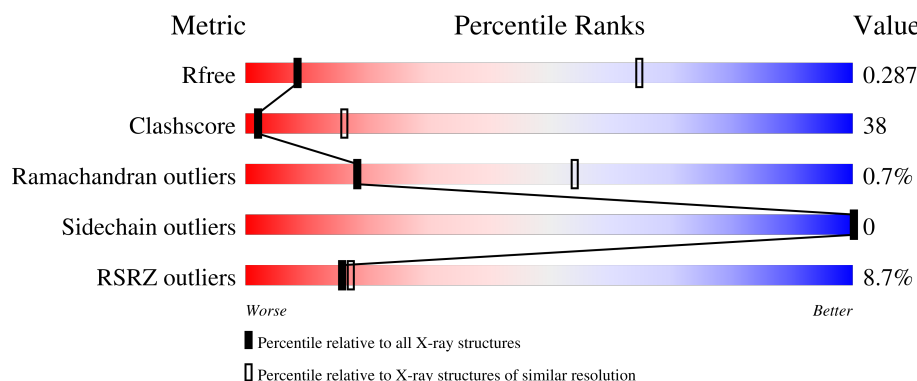
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1082 (4.20-3.80)
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)
RSRZ outliers	180081	1082 (4.20-3.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	376	7% 32% 35% 34%
1	B	376	3% 32% 33% 34%
1	C	376	8% 32% 33% 34%
1	D	376	5% 30% 35% 34%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	D	403	-	-	X	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15472 atoms, of which 7776 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Aquaporin-7.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	249	Total	C	H	N	O	S	0	0	0
			3826	1266	1920	308	319	13			
1	C	249	Total	C	H	N	O	S	0	0	0
			3826	1266	1920	308	319	13			
1	B	249	Total	C	H	N	O	S	0	0	0
			3826	1266	1920	308	319	13			
1	D	249	Total	C	H	N	O	S	0	0	0
			3826	1266	1920	308	319	13			

There are 136 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	343	GLY	-	expression tag	UNP O14520
A	344	GLY	-	expression tag	UNP O14520
A	345	SER	-	expression tag	UNP O14520
A	346	LEU	-	expression tag	UNP O14520
A	347	GLU	-	expression tag	UNP O14520
A	348	VAL	-	expression tag	UNP O14520
A	349	LEU	-	expression tag	UNP O14520
A	350	PHE	-	expression tag	UNP O14520
A	351	GLN	-	expression tag	UNP O14520
A	352	GLY	-	expression tag	UNP O14520
A	353	PRO	-	expression tag	UNP O14520
A	354	ALA	-	expression tag	UNP O14520
A	355	ALA	-	expression tag	UNP O14520
A	356	TYR	-	expression tag	UNP O14520
A	357	PRO	-	expression tag	UNP O14520
A	358	TYR	-	expression tag	UNP O14520
A	359	ASP	-	expression tag	UNP O14520
A	360	VAL	-	expression tag	UNP O14520
A	361	PRO	-	expression tag	UNP O14520
A	362	ASP	-	expression tag	UNP O14520
A	363	TYR	-	expression tag	UNP O14520

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Chain	Residue	Modelled	Actual	Comment	Reference
A	364	ALA	-	expression tag	UNP O14520
A	365	ALA	-	expression tag	UNP O14520
A	366	ALA	-	expression tag	UNP O14520
A	367	HIS	-	expression tag	UNP O14520
A	368	HIS	-	expression tag	UNP O14520
A	369	HIS	-	expression tag	UNP O14520
A	370	HIS	-	expression tag	UNP O14520
A	371	HIS	-	expression tag	UNP O14520
A	372	HIS	-	expression tag	UNP O14520
A	373	HIS	-	expression tag	UNP O14520
A	374	HIS	-	expression tag	UNP O14520
A	375	HIS	-	expression tag	UNP O14520
A	376	HIS	-	expression tag	UNP O14520
C	343	GLY	-	expression tag	UNP O14520
C	344	GLY	-	expression tag	UNP O14520
C	345	SER	-	expression tag	UNP O14520
C	346	LEU	-	expression tag	UNP O14520
C	347	GLU	-	expression tag	UNP O14520
C	348	VAL	-	expression tag	UNP O14520
C	349	LEU	-	expression tag	UNP O14520
C	350	PHE	-	expression tag	UNP O14520
C	351	GLN	-	expression tag	UNP O14520
C	352	GLY	-	expression tag	UNP O14520
C	353	PRO	-	expression tag	UNP O14520
C	354	ALA	-	expression tag	UNP O14520
C	355	ALA	-	expression tag	UNP O14520
C	356	TYR	-	expression tag	UNP O14520
C	357	PRO	-	expression tag	UNP O14520
C	358	TYR	-	expression tag	UNP O14520
C	359	ASP	-	expression tag	UNP O14520
C	360	VAL	-	expression tag	UNP O14520
C	361	PRO	-	expression tag	UNP O14520
C	362	ASP	-	expression tag	UNP O14520
C	363	TYR	-	expression tag	UNP O14520
C	364	ALA	-	expression tag	UNP O14520
C	365	ALA	-	expression tag	UNP O14520
C	366	ALA	-	expression tag	UNP O14520
C	367	HIS	-	expression tag	UNP O14520
C	368	HIS	-	expression tag	UNP O14520
C	369	HIS	-	expression tag	UNP O14520
C	370	HIS	-	expression tag	UNP O14520
C	371	HIS	-	expression tag	UNP O14520

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Chain	Residue	Modelled	Actual	Comment	Reference
C	372	HIS	-	expression tag	UNP O14520
C	373	HIS	-	expression tag	UNP O14520
C	374	HIS	-	expression tag	UNP O14520
C	375	HIS	-	expression tag	UNP O14520
C	376	HIS	-	expression tag	UNP O14520
B	343	GLY	-	expression tag	UNP O14520
B	344	GLY	-	expression tag	UNP O14520
B	345	SER	-	expression tag	UNP O14520
B	346	LEU	-	expression tag	UNP O14520
B	347	GLU	-	expression tag	UNP O14520
B	348	VAL	-	expression tag	UNP O14520
B	349	LEU	-	expression tag	UNP O14520
B	350	PHE	-	expression tag	UNP O14520
B	351	GLN	-	expression tag	UNP O14520
B	352	GLY	-	expression tag	UNP O14520
B	353	PRO	-	expression tag	UNP O14520
B	354	ALA	-	expression tag	UNP O14520
B	355	ALA	-	expression tag	UNP O14520
B	356	TYR	-	expression tag	UNP O14520
B	357	PRO	-	expression tag	UNP O14520
B	358	TYR	-	expression tag	UNP O14520
B	359	ASP	-	expression tag	UNP O14520
B	360	VAL	-	expression tag	UNP O14520
B	361	PRO	-	expression tag	UNP O14520
B	362	ASP	-	expression tag	UNP O14520
B	363	TYR	-	expression tag	UNP O14520
B	364	ALA	-	expression tag	UNP O14520
B	365	ALA	-	expression tag	UNP O14520
B	366	ALA	-	expression tag	UNP O14520
B	367	HIS	-	expression tag	UNP O14520
B	368	HIS	-	expression tag	UNP O14520
B	369	HIS	-	expression tag	UNP O14520
B	370	HIS	-	expression tag	UNP O14520
B	371	HIS	-	expression tag	UNP O14520
B	372	HIS	-	expression tag	UNP O14520
B	373	HIS	-	expression tag	UNP O14520
B	374	HIS	-	expression tag	UNP O14520
B	375	HIS	-	expression tag	UNP O14520
B	376	HIS	-	expression tag	UNP O14520
D	343	GLY	-	expression tag	UNP O14520
D	344	GLY	-	expression tag	UNP O14520
D	345	SER	-	expression tag	UNP O14520

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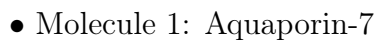
Chain	Residue	Modelled	Actual	Comment	Reference
D	346	LEU	-	expression tag	UNP O14520
D	347	GLU	-	expression tag	UNP O14520
D	348	VAL	-	expression tag	UNP O14520
D	349	LEU	-	expression tag	UNP O14520
D	350	PHE	-	expression tag	UNP O14520
D	351	GLN	-	expression tag	UNP O14520
D	352	GLY	-	expression tag	UNP O14520
D	353	PRO	-	expression tag	UNP O14520
D	354	ALA	-	expression tag	UNP O14520
D	355	ALA	-	expression tag	UNP O14520
D	356	TYR	-	expression tag	UNP O14520
D	357	PRO	-	expression tag	UNP O14520
D	358	TYR	-	expression tag	UNP O14520
D	359	ASP	-	expression tag	UNP O14520
D	360	VAL	-	expression tag	UNP O14520
D	361	PRO	-	expression tag	UNP O14520
D	362	ASP	-	expression tag	UNP O14520
D	363	TYR	-	expression tag	UNP O14520
D	364	ALA	-	expression tag	UNP O14520
D	365	ALA	-	expression tag	UNP O14520
D	366	ALA	-	expression tag	UNP O14520
D	367	HIS	-	expression tag	UNP O14520
D	368	HIS	-	expression tag	UNP O14520
D	369	HIS	-	expression tag	UNP O14520
D	370	HIS	-	expression tag	UNP O14520
D	371	HIS	-	expression tag	UNP O14520
D	372	HIS	-	expression tag	UNP O14520
D	373	HIS	-	expression tag	UNP O14520
D	374	HIS	-	expression tag	UNP O14520
D	375	HIS	-	expression tag	UNP O14520
D	376	HIS	-	expression tag	UNP O14520

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			14	3	8	3		
2	A	1	Total	C	H	O	0	0
			14	3	8	3		
2	C	1	Total	C	H	O	0	0
			14	3	8	3		
2	C	1	Total	C	H	O	0	0
			14	3	8	3		
2	C	1	Total	C	H	O	0	0
			14	3	8	3		
2	B	1	Total	C	H	O	0	0
			14	3	8	3		
2	B	1	Total	C	H	O	0	0
			14	3	8	3		
2	B	1	Total	C	H	O	14	0
			14	3	8	3		
2	D	1	Total	C	H	O	0	0
			14	3	8	3		
2	D	1	Total	C	H	O	0	0
			14	3	8	3		
2	D	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 1: Aquaporin-7



[illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	F 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	385.66Å 385.66Å 385.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.71 – 4.00 32.71 – 4.00	Depositor EDS
% Data completeness (in resolution range)	87.9 (32.71-4.00) 93.1 (32.71-4.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	50.00 (at 4.00Å)	Xtrriage
Refinement program	PHENIX (1.14_3235: ???)	Depositor
R, R_{free}	0.232 , 0.277 0.243 , 0.287	Depositor DCC
R_{free} test set	599 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å ²)	132.5	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.75 , 709.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	15472	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 88.55 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.7106e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.28	0/1963	0.56	0/2676
1	B	0.28	0/1963	0.59	1/2676 (0.0%)
1	C	0.29	0/1963	0.56	0/2676
1	D	0.30	0/1963	0.59	1/2676 (0.0%)
All	All	0.29	0/7852	0.57	2/10704 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	253	TRP	N-CA-C	6.71	125.09	110.80
1	B	253	TRP	N-CA-C	6.04	123.66	110.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	194	ASN	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1906	1920	1920	160	1
1	B	1906	1920	1920	167	1
1	C	1906	1920	1920	151	1
1	D	1906	1920	1920	171	1
2	A	12	16	16	0	0
2	B	18	24	24	0	0
2	C	18	24	24	3	0
2	D	24	32	32	7	0
All	All	7696	7776	7776	585	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 38.

All (585) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:MET:HE1	1:A:111:LYS:HG2	1.39	1.05
1:D:40:GLU:HA	1:D:93:MET:HE1	1.49	0.93
1:D:45:TYR:OH	1:D:129:THR:OG1	1.87	0.93
1:B:162:LEU:HD13	1:B:253:TRP:CH2	2.04	0.92
1:B:40:GLU:HA	1:B:93:MET:HE1	1.50	0.92
1:D:60:LEU:CD1	1:D:138:ILE:HD13	2.00	0.92
1:A:39:ALA:HB3	1:A:88:ILE:HD11	1.51	0.91
1:A:40:GLU:HA	1:A:93:MET:HE1	1.53	0.90
1:B:173:LEU:CD2	1:D:133:LEU:HD21	2.02	0.88
1:D:182:LEU:HD13	1:D:225:ILE:HG21	1.56	0.88
1:C:178:LEU:HD21	1:C:217:LEU:HD22	1.55	0.87
1:C:45:TYR:OH	1:C:129:THR:OG1	1.92	0.87
1:C:152:VAL:HG12	1:B:146:LEU:HD22	1.57	0.86
1:A:162:LEU:HD13	1:A:253:TRP:CH2	2.12	0.85
1:C:178:LEU:HD12	1:C:213:ILE:HG23	1.57	0.85
1:B:173:LEU:HD22	1:D:133:LEU:HD21	1.56	0.84
1:D:162:LEU:HD13	1:D:253:TRP:CH2	2.12	0.83
1:B:182:LEU:HD13	1:B:225:ILE:HG21	1.60	0.83
1:B:205:VAL:HG23	1:D:80:MET:HE1	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:SER:OG	1:C:64:TYR:O	1.97	0.82
1:C:178:LEU:HD11	1:C:217:LEU:HD13	1.61	0.82
1:A:131:TYR:HE1	1:A:138:ILE:HD13	1.45	0.81
1:C:32:LYS:HG3	1:C:35:ARG:HG3	1.62	0.80
1:B:237:THR:HB	1:B:246:VAL:HG21	1.63	0.80
1:D:59:VAL:HG13	1:D:60:LEU:HG	1.64	0.80
1:D:195:ASN:HB3	1:D:196:PRO:HD3	1.65	0.79
1:B:173:LEU:HD22	1:D:133:LEU:CD2	2.12	0.79
1:A:146:LEU:HB3	1:D:152:VAL:HG13	1.67	0.77
1:D:60:LEU:HD13	1:D:138:ILE:HD13	1.67	0.77
1:A:237:THR:HB	1:A:246:VAL:HG21	1.67	0.77
1:D:274:VAL:HG13	1:D:275:PHE:CD1	2.20	0.76
1:C:39:ALA:CB	1:C:88:ILE:HD11	2.16	0.76
1:A:170:ARG:HH12	1:C:136:THR:HG23	1.50	0.76
1:A:270:ILE:O	1:A:274:VAL:HG12	1.86	0.76
1:A:205:VAL:HG23	1:C:80:MET:HE1	1.66	0.75
1:D:45:TYR:O	1:D:49:VAL:HG12	1.85	0.75
1:B:162:LEU:HD11	1:B:166:MET:SD	2.25	0.75
1:D:237:THR:HB	1:D:246:VAL:HG21	1.67	0.75
1:A:167:THR:HG21	1:A:169:TRP:CZ2	2.22	0.75
1:C:57:HIS:CE1	1:C:63:LYS:HB2	2.23	0.74
1:C:237:THR:HB	1:C:246:VAL:HG21	1.70	0.74
1:D:186:LEU:HD11	2:D:403:GOL:C3	2.18	0.74
1:D:54:SER:O	1:D:63:LYS:NZ	2.17	0.74
1:D:57:HIS:CE1	1:D:63:LYS:HB3	2.23	0.74
1:D:91:ALA:O	2:D:403:GOL:O3	2.03	0.74
1:B:167:THR:HG21	1:B:169:TRP:CZ2	2.23	0.73
1:C:184:LEU:HD12	1:C:271:ILE:HG23	1.70	0.73
1:B:274:VAL:HG13	1:B:275:PHE:CD1	2.23	0.73
1:B:198:LEU:HD13	1:B:201:THR:HB	1.69	0.73
1:C:162:LEU:HD13	1:C:253:TRP:CH2	2.24	0.73
1:A:104:LEU:HD13	1:A:104:LEU:O	1.87	0.73
1:A:36:GLU:HA	1:A:88:ILE:HD13	1.71	0.72
1:C:59:VAL:HG23	1:C:60:LEU:HG	1.72	0.72
1:B:198:LEU:HD13	1:B:201:THR:C	2.14	0.72
1:D:167:THR:HG21	1:D:169:TRP:CZ2	2.25	0.72
1:A:57:HIS:CE1	1:A:63:LYS:HB2	2.24	0.71
1:B:57:HIS:CE1	1:B:63:LYS:HB2	2.25	0.71
1:B:74:PHE:O	1:B:78:VAL:HG23	1.89	0.71
1:D:195:ASN:HB3	1:D:196:PRO:CD	2.20	0.71
1:B:45:TYR:CZ	1:B:126:ALA:HA	2.25	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLN:O	1:A:228:SER:OG	2.08	0.71
1:A:45:TYR:OH	1:A:129:THR:OG1	2.06	0.71
1:B:270:ILE:O	1:B:274:VAL:HG12	1.91	0.71
1:C:167:THR:HG21	1:C:169:TRP:CZ2	2.26	0.70
1:C:33:MET:HE1	1:C:115:TYR:CE1	2.27	0.70
1:D:45:TYR:CZ	1:D:126:ALA:HA	2.27	0.70
1:B:189:ILE:HA	1:D:83:HIS:NE2	2.07	0.70
1:B:198:LEU:CD1	1:B:201:THR:C	2.65	0.69
1:C:167:THR:HG21	1:C:169:TRP:CH2	2.28	0.68
1:C:195:ASN:HB3	1:C:196:PRO:HD3	1.75	0.68
1:C:40:GLU:HA	1:C:93:MET:HE1	1.75	0.68
1:C:45:TYR:O	1:C:49:VAL:HG12	1.94	0.67
1:C:52:LEU:HD22	1:C:130:ILE:HG21	1.75	0.67
1:A:67:TYR:CE2	1:A:215:VAL:HG22	2.28	0.67
1:B:198:LEU:HD12	1:B:202:GLU:N	2.09	0.67
1:A:52:LEU:O	1:A:52:LEU:HD23	1.93	0.67
1:A:52:LEU:HD22	1:A:130:ILE:HG21	1.76	0.67
1:C:131:TYR:CE1	1:C:138:ILE:HD13	2.29	0.67
1:A:39:ALA:HB3	1:A:88:ILE:CD1	2.25	0.67
1:A:59:VAL:HG13	1:A:60:LEU:HG	1.77	0.67
1:A:162:LEU:HD11	1:A:166:MET:SD	2.36	0.66
1:B:45:TYR:OH	1:B:129:THR:OG1	2.11	0.66
1:A:47:MET:HG2	1:A:78:VAL:HG12	1.77	0.66
1:A:150:GLY:N	1:A:151:PRO:HD3	2.10	0.66
1:A:165:HIS:O	1:C:136:THR:OG1	2.11	0.66
1:A:274:VAL:HG13	1:A:275:PHE:CD1	2.31	0.66
1:B:173:LEU:HD23	1:D:133:LEU:HD21	1.77	0.65
1:C:45:TYR:CZ	1:C:126:ALA:HA	2.31	0.65
1:B:52:LEU:HD23	1:B:52:LEU:O	1.95	0.65
1:A:45:TYR:CZ	1:A:126:ALA:HA	2.32	0.64
1:D:276:ILE:O	1:D:279:THR:OG1	2.10	0.64
1:A:138:ILE:HG13	1:A:139:LEU:N	2.11	0.64
1:B:44:THR:N	1:B:93:MET:SD	2.71	0.64
1:B:33:MET:HE1	1:B:115:TYR:CE1	2.32	0.64
1:B:205:VAL:CG2	1:D:80:MET:HE1	2.27	0.64
1:D:57:HIS:O	1:D:63:LYS:HG2	1.98	0.63
1:A:146:LEU:HB3	1:D:152:VAL:CG1	2.27	0.63
1:C:152:VAL:HA	1:B:146:LEU:HD13	1.79	0.63
1:C:152:VAL:CG1	1:B:146:LEU:HD22	2.27	0.63
1:C:150:GLY:N	1:C:151:PRO:HD3	2.14	0.62
1:B:198:LEU:CD1	1:B:202:GLU:N	2.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:LEU:HD11	1:B:202:GLU:HA	1.81	0.62
1:A:48:MET:O	1:A:52:LEU:N	2.20	0.62
1:A:88:ILE:HD12	1:A:89:SER:N	2.14	0.62
1:C:39:ALA:HB3	1:C:88:ILE:HD11	1.80	0.62
1:C:45:TYR:HH	1:C:129:THR:HG1	1.43	0.62
1:D:248:SER:O	1:D:251:GLU:O	2.18	0.62
1:D:52:LEU:HD21	1:D:158:PHE:CD1	2.35	0.61
1:D:44:THR:N	1:D:93:MET:SD	2.73	0.61
1:C:39:ALA:HB2	1:C:88:ILE:HD11	1.83	0.61
1:D:33:MET:HE1	1:D:115:TYR:CE1	2.36	0.61
1:A:39:ALA:CB	1:A:88:ILE:HD11	2.28	0.60
1:C:52:LEU:O	1:C:52:LEU:HD23	2.01	0.60
1:D:82:VAL:HG22	2:D:403:GOL:H12	1.82	0.60
1:B:252:ASN:O	1:B:254:TRP:N	2.31	0.60
1:A:256:VAL:HG13	1:A:257:PRO:HD3	1.82	0.60
1:A:279:THR:HB	1:C:35:ARG:HD2	1.83	0.60
1:C:32:LYS:CG	1:C:35:ARG:HG3	2.29	0.60
1:C:33:MET:HE1	1:C:115:TYR:CD1	2.37	0.60
1:D:167:THR:HG21	1:D:169:TRP:CH2	2.36	0.60
1:D:48:MET:O	1:D:52:LEU:N	2.20	0.60
1:B:32:LYS:HB3	1:B:35:ARG:NE	2.16	0.60
1:B:104:LEU:O	1:B:104:LEU:HG	2.01	0.60
1:D:274:VAL:HG13	1:D:275:PHE:HD1	1.64	0.59
1:A:131:TYR:CE1	1:A:138:ILE:HD13	2.32	0.59
1:C:111:LYS:HD3	1:C:115:TYR:OH	2.02	0.59
1:D:60:LEU:HD11	1:D:138:ILE:HD13	1.82	0.59
1:B:60:LEU:HD23	1:B:137:ALA:O	2.02	0.58
1:A:194:ASN:OD1	1:C:87:ARG:HB2	2.03	0.58
1:B:45:TYR:O	1:B:49:VAL:HG12	2.04	0.58
1:A:94:ASN:HB3	1:A:97:VAL:HG22	1.86	0.58
1:D:44:THR:OG1	1:D:93:MET:SD	2.58	0.58
1:D:198:LEU:O	1:D:202:GLU:HG3	2.04	0.58
1:A:146:LEU:O	1:D:152:VAL:HG21	2.04	0.58
1:A:219:MET:HB2	1:C:57:HIS:CE1	2.38	0.58
1:C:39:ALA:HA	1:C:42:MET:HE3	1.86	0.58
1:B:178:LEU:HD21	1:B:217:LEU:HD22	1.85	0.58
1:A:152:VAL:HG11	1:D:146:LEU:HB3	1.86	0.58
1:A:279:THR:HA	1:C:35:ARG:HD3	1.86	0.58
1:D:92:HIS:O	2:D:403:GOL:O2	2.21	0.58
1:B:67:TYR:CE1	1:D:63:LYS:HD2	2.39	0.57
1:C:195:ASN:HB3	1:C:196:PRO:CD	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:256:VAL:CG1	1:B:257:PRO:HD3	2.35	0.57
1:B:189:ILE:HG22	1:B:202:GLU:HG3	1.87	0.57
1:D:52:LEU:HG	1:D:130:ILE:HG21	1.87	0.57
1:A:198:LEU:O	1:A:202:GLU:HG3	2.04	0.57
1:C:52:LEU:HD22	1:C:130:ILE:CG2	2.34	0.57
1:A:88:ILE:HD12	1:A:89:SER:HB3	1.86	0.57
1:C:135:TYR:O	1:C:138:ILE:HG12	2.04	0.57
1:C:219:MET:O	1:C:219:MET:HG3	2.04	0.57
1:B:39:ALA:HB1	1:B:85:ALA:HB1	1.87	0.57
1:B:195:ASN:CG	1:B:196:PRO:HD3	2.30	0.57
1:D:182:LEU:HD13	1:D:225:ILE:CG2	2.32	0.56
1:A:150:GLY:N	1:A:151:PRO:CD	2.68	0.56
1:B:98:THR:HG22	1:B:112:PHE:HD1	1.70	0.56
1:B:226:ASN:HB3	1:B:229:ARG:HG2	1.86	0.56
1:D:52:LEU:HD21	1:D:158:PHE:HD1	1.70	0.56
1:D:131:TYR:CE1	1:D:138:ILE:HG13	2.41	0.56
1:C:113:PRO:O	1:C:116:VAL:HG12	2.05	0.56
1:B:198:LEU:CD1	1:B:202:GLU:HA	2.36	0.56
1:C:168:LEU:HD22	1:C:252:ASN:HD21	1.71	0.56
1:A:39:ALA:HA	1:A:42:MET:HE3	1.88	0.56
1:A:152:VAL:CG1	1:D:146:LEU:HB3	2.36	0.56
1:C:72:LEU:HA	1:C:211:VAL:HG22	1.88	0.56
1:B:194:ASN:OD1	1:D:87:ARG:HB2	2.05	0.55
1:D:162:LEU:HD11	1:D:166:MET:SD	2.45	0.55
1:D:205:VAL:HA	1:D:208:ILE:HG12	1.87	0.55
1:A:44:THR:OG1	1:A:119:GLN:HB2	2.06	0.55
1:C:48:MET:HB3	1:C:126:ALA:HB1	1.87	0.55
1:B:189:ILE:HD13	1:B:205:VAL:HG13	1.88	0.55
1:B:32:LYS:HB3	1:B:35:ARG:CD	2.36	0.55
1:B:94:ASN:HB3	1:B:97:VAL:HB	1.88	0.55
1:C:146:LEU:HB2	1:B:152:VAL:CG1	2.36	0.55
1:B:93:MET:HE2	1:B:93:MET:H	1.70	0.55
1:A:44:THR:N	1:A:93:MET:SD	2.79	0.55
1:B:276:ILE:HD11	1:D:42:MET:SD	2.47	0.55
1:C:150:GLY:N	1:C:151:PRO:CD	2.70	0.55
1:C:205:VAL:HA	1:C:208:ILE:HG12	1.89	0.55
1:D:74:PHE:O	1:D:78:VAL:HG23	2.06	0.55
1:A:189:ILE:HD13	1:A:205:VAL:HG13	1.89	0.55
1:A:162:LEU:HD13	1:A:253:TRP:CZ2	2.42	0.54
1:B:198:LEU:HD12	1:B:198:LEU:O	2.07	0.54
1:A:57:HIS:ND1	1:A:63:LYS:HB2	2.21	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:THR:OG1	1:B:93:MET:SD	2.58	0.54
1:D:192:GLN:NE2	1:D:193:GLU:OE2	2.40	0.54
1:A:205:VAL:CG2	1:C:80:MET:HE1	2.37	0.54
1:D:60:LEU:CD1	1:D:138:ILE:CD1	2.83	0.54
1:B:67:TYR:HE1	1:D:65:GLY:CA	2.21	0.54
1:D:149:THR:O	1:D:149:THR:HG23	2.07	0.54
1:B:178:LEU:HD11	1:B:217:LEU:CB	2.37	0.54
1:B:212:ILE:HD11	1:D:76:PHE:HB3	1.89	0.54
1:D:48:MET:HB3	1:D:126:ALA:HB1	1.89	0.54
1:D:105:GLY:O	1:D:106:ARG:HG2	2.08	0.54
1:C:47:MET:HG2	1:C:78:VAL:HG22	1.90	0.54
1:C:174:ASN:O	1:C:178:LEU:HD23	2.08	0.54
1:C:184:LEU:HD12	1:C:271:ILE:CG2	2.38	0.53
1:A:195:ASN:OD1	1:A:196:PRO:HD2	2.07	0.53
1:B:219:MET:HB2	1:D:57:HIS:CE1	2.43	0.53
1:D:32:LYS:HG3	1:D:35:ARG:HG3	1.89	0.53
1:A:37:PHE:CZ	1:A:118:GLY:HA2	2.43	0.53
1:A:195:ASN:CG	1:A:196:PRO:CD	2.82	0.53
1:B:230:ASP:O	1:B:234:ARG:HG3	2.08	0.53
1:D:255:TRP:O	1:D:258:VAL:HG12	2.08	0.53
1:D:32:LYS:HG2	1:D:35:ARG:NE	2.22	0.53
1:C:104:LEU:HG	1:C:104:LEU:O	2.08	0.53
1:C:117:LEU:O	1:C:117:LEU:HD23	2.08	0.53
1:B:52:LEU:HD13	1:B:130:ILE:HG21	1.89	0.53
1:C:206:ILE:HG22	2:C:403:GOL:O2	2.09	0.53
1:D:178:LEU:HD21	1:D:217:LEU:HD13	1.89	0.53
1:A:48:MET:HB3	1:A:126:ALA:HB1	1.91	0.52
1:B:150:GLY:N	1:B:151:PRO:CD	2.72	0.52
1:D:195:ASN:CB	1:D:196:PRO:HD3	2.38	0.52
1:A:189:ILE:HA	1:C:83:HIS:NE2	2.24	0.52
1:B:279:THR:HG22	1:D:35:ARG:HG2	1.90	0.52
1:A:230:ASP:O	1:A:234:ARG:HG3	2.09	0.52
1:C:194:ASN:O	1:C:195:ASN:C	2.51	0.52
1:B:57:HIS:ND1	1:B:63:LYS:HB2	2.24	0.52
1:B:142:SER:OG	1:B:147:MET:O	2.28	0.52
1:A:102:CYS:HA	1:A:107:VAL:HG12	1.92	0.52
1:B:52:LEU:HD11	1:B:158:PHE:HD1	1.74	0.52
1:D:98:THR:HG22	1:D:112:PHE:HD1	1.74	0.52
1:D:277:GLY:O	1:D:280:ILE:HG12	2.10	0.52
1:C:94:ASN:CG	1:C:226:ASN:OD1	2.53	0.52
1:C:178:LEU:HD12	1:C:213:ILE:CG2	2.33	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:LEU:HD11	1:B:158:PHE:CD1	2.45	0.52
1:B:237:THR:HG23	1:B:242:TRP:CE3	2.44	0.52
1:D:36:GLU:HG2	1:D:89:SER:HA	1.92	0.52
1:C:198:LEU:O	1:C:202:GLU:HG3	2.10	0.52
1:B:67:TYR:CE2	1:B:215:VAL:HG22	2.45	0.52
1:D:249:ASN:O	1:D:249:ASN:ND2	2.43	0.52
1:A:93:MET:O	1:A:226:ASN:ND2	2.43	0.51
1:D:149:THR:C	1:D:151:PRO:CD	2.83	0.51
1:A:52:LEU:HD22	1:A:130:ILE:CG2	2.39	0.51
1:C:149:THR:HG23	1:C:149:THR:O	2.08	0.51
1:C:162:LEU:HD11	1:C:166:MET:SD	2.50	0.51
1:B:149:THR:O	1:B:149:THR:HG23	2.11	0.51
1:A:72:LEU:HA	1:A:211:VAL:HG22	1.93	0.51
1:B:35:ARG:HG3	1:B:88:ILE:HD11	1.91	0.51
1:B:198:LEU:CD1	1:B:202:GLU:CA	2.88	0.51
1:D:72:LEU:HA	1:D:211:VAL:HG22	1.93	0.51
1:B:59:VAL:HB	1:B:60:LEU:HD12	1.93	0.51
1:B:105:GLY:O	1:B:106:ARG:HG2	2.11	0.51
1:D:93:MET:O	1:D:226:ASN:ND2	2.44	0.51
1:D:150:GLY:N	1:D:151:PRO:CD	2.73	0.51
1:A:68:LEU:HD22	1:C:69:GLY:HA2	1.93	0.51
1:C:119:GLN:O	1:C:228:SER:OG	2.28	0.51
1:C:255:TRP:O	1:C:258:VAL:HG12	2.10	0.51
1:B:149:THR:OG1	1:B:151:PRO:CD	2.58	0.51
1:B:182:LEU:HA	1:B:209:LEU:HD21	1.92	0.51
1:D:39:ALA:HA	1:D:42:MET:HE3	1.92	0.51
1:D:60:LEU:HD11	1:D:138:ILE:CD1	2.41	0.51
1:C:198:LEU:HD22	1:C:201:THR:OG1	2.11	0.51
1:B:193:GLU:O	1:D:87:ARG:NH1	2.44	0.51
1:A:32:LYS:HB3	1:A:35:ARG:NE	2.26	0.50
1:B:32:LYS:NZ	1:B:35:ARG:HB3	2.26	0.50
1:D:33:MET:SD	1:D:115:TYR:HE1	2.34	0.50
1:D:186:LEU:HD11	2:D:403:GOL:H31	1.92	0.50
1:A:182:LEU:HA	1:A:209:LEU:HD21	1.93	0.50
1:A:255:TRP:O	1:A:258:VAL:HG12	2.12	0.50
1:C:48:MET:O	1:C:52:LEU:N	2.25	0.50
1:B:33:MET:HE3	1:B:114:VAL:CG1	2.40	0.50
1:D:40:GLU:CA	1:D:93:MET:HE1	2.30	0.50
1:A:149:THR:OG1	1:A:151:PRO:CD	2.60	0.50
1:D:57:HIS:CG	1:D:63:LYS:HD3	2.47	0.50
1:D:198:LEU:O	1:D:199:PRO:C	2.55	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:THR:HG22	1:C:112:PHE:HD1	1.76	0.50
1:B:60:LEU:HB3	1:B:137:ALA:HB1	1.94	0.50
1:D:149:THR:OG1	1:D:151:PRO:CD	2.60	0.50
1:D:161:TYR:CG	1:D:249:ASN:OD1	2.65	0.50
1:B:72:LEU:HA	1:B:211:VAL:HG22	1.93	0.50
1:B:216:SER:OG	1:B:217:LEU:HD12	2.11	0.50
1:B:66:SER:HB3	1:D:64:TYR:O	2.12	0.50
1:B:167:THR:HG21	1:B:169:TRP:CH2	2.45	0.50
1:D:111:LYS:HD3	1:D:115:TYR:OH	2.12	0.50
1:C:111:LYS:HD3	1:C:115:TYR:CZ	2.47	0.50
1:C:249:ASN:ND2	1:C:249:ASN:O	2.45	0.49
1:C:192:GLN:NE2	1:C:193:GLU:OE2	2.46	0.49
1:B:178:LEU:HD21	1:B:217:LEU:HD13	1.95	0.49
1:B:253:TRP:CZ3	1:B:256:VAL:HG11	2.47	0.49
1:A:152:VAL:HG21	1:D:146:LEU:O	2.12	0.49
1:A:178:LEU:HD21	1:A:217:LEU:HD22	1.93	0.49
1:C:160:THR:HG21	1:C:224:ALA:HB2	1.95	0.49
1:C:115:TYR:O	1:C:119:GLN:HG2	2.12	0.49
1:C:149:THR:OG1	1:C:151:PRO:HD2	2.12	0.49
1:D:154:THR:O	1:D:157:ILE:HG12	2.13	0.49
1:A:237:THR:HG23	1:A:242:TRP:CE3	2.48	0.49
1:A:279:THR:HB	1:C:35:ARG:CD	2.42	0.49
1:C:131:TYR:HE1	1:C:138:ILE:HD13	1.78	0.49
1:C:256:VAL:CG1	1:C:257:PRO:HD3	2.42	0.49
1:B:149:THR:C	1:B:151:PRO:CD	2.85	0.49
1:C:74:PHE:O	1:C:78:VAL:HG23	2.12	0.48
1:C:138:ILE:CG1	1:C:139:LEU:N	2.75	0.48
1:B:194:ASN:O	1:B:195:ASN:O	2.31	0.48
1:D:94:ASN:HB3	1:D:97:VAL:HG22	1.94	0.48
1:A:108:PRO:HG3	1:A:111:LYS:HE2	1.96	0.48
1:D:111:LYS:HB3	1:D:115:TYR:CE1	2.49	0.48
1:A:32:LYS:HG2	1:A:35:ARG:HB2	1.96	0.48
1:A:72:LEU:HD11	1:A:76:PHE:CZ	2.48	0.48
1:A:219:MET:CB	1:C:57:HIS:CE1	2.96	0.48
1:B:277:GLY:O	1:B:280:ILE:HG12	2.14	0.48
1:D:113:PRO:O	1:D:116:VAL:HG12	2.13	0.48
1:D:142:SER:OG	1:D:147:MET:O	2.32	0.48
1:D:178:LEU:HD11	1:D:217:LEU:CB	2.44	0.48
1:D:168:LEU:CD1	1:D:252:ASN:HD22	2.27	0.48
1:A:149:THR:OG1	1:A:151:PRO:HD2	2.14	0.48
1:A:197:ALA:O	1:A:199:PRO:HD3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:GLU:CA	1:B:93:MET:HE1	2.33	0.48
1:D:52:LEU:CG	1:D:130:ILE:HG21	2.43	0.48
1:A:93:MET:H	1:A:93:MET:HE2	1.79	0.48
1:B:52:LEU:HD22	1:B:130:ILE:HG21	1.96	0.48
1:B:57:HIS:HD1	1:B:61:ASN:HB2	1.79	0.48
1:D:52:LEU:CD2	1:D:130:ILE:HG21	2.44	0.48
1:A:44:THR:OG1	1:A:93:MET:SD	2.59	0.47
1:C:111:LYS:HD3	1:C:115:TYR:CE1	2.49	0.47
1:C:192:GLN:OE1	1:C:192:GLN:N	2.31	0.47
1:B:183:GLN:NE2	1:B:187:PHE:CZ	2.82	0.47
1:A:149:THR:HG23	1:A:149:THR:O	2.13	0.47
1:A:160:THR:HB	1:A:221:THR:O	2.14	0.47
1:B:37:PHE:HD2	1:B:114:VAL:HG12	1.80	0.47
1:B:38:LEU:O	1:B:38:LEU:HD23	2.14	0.47
1:A:117:LEU:HD23	1:A:117:LEU:O	2.14	0.47
1:C:111:LYS:O	1:C:115:TYR:CD1	2.67	0.47
1:B:195:ASN:CG	1:B:196:PRO:CD	2.87	0.47
1:B:67:TYR:HE1	1:D:65:GLY:HA2	1.80	0.47
1:B:39:ALA:HA	1:B:42:MET:HE3	1.97	0.47
1:B:198:LEU:HD12	1:B:202:GLU:CA	2.45	0.47
1:B:267:LEU:HD22	1:B:270:ILE:HD11	1.97	0.47
1:A:36:GLU:OE1	1:A:115:TYR:OH	2.33	0.47
1:A:195:ASN:CG	1:A:196:PRO:HD2	2.40	0.47
1:C:138:ILE:HG13	1:C:139:LEU:N	2.30	0.47
1:B:67:TYR:CE1	1:D:65:GLY:CA	2.98	0.47
1:D:32:LYS:CG	1:D:35:ARG:HG3	2.45	0.47
1:D:194:ASN:O	1:D:195:ASN:C	2.57	0.47
1:D:237:THR:HG23	1:D:242:TRP:CE3	2.50	0.47
1:D:256:VAL:CG1	1:D:257:PRO:HD3	2.44	0.47
1:A:256:VAL:CG1	1:A:257:PRO:HD3	2.45	0.47
1:D:32:LYS:N	1:D:35:ARG:HD2	2.29	0.47
1:A:149:THR:C	1:A:151:PRO:CD	2.88	0.47
1:C:162:LEU:HD13	1:C:253:TRP:CZ2	2.49	0.47
1:C:105:GLY:O	1:C:106:ARG:HG2	2.15	0.46
1:D:93:MET:HE2	1:D:93:MET:H	1.78	0.46
1:C:146:LEU:HB2	1:B:152:VAL:HG13	1.98	0.46
1:B:72:LEU:HD11	1:B:76:PHE:CZ	2.50	0.46
1:C:253:TRP:CH2	1:C:256:VAL:HG11	2.50	0.46
1:A:178:LEU:HD11	1:A:217:LEU:CB	2.45	0.46
1:A:251:GLU:O	1:A:252:ASN:C	2.59	0.46
1:B:195:ASN:OD1	1:B:196:PRO:HD3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:SER:O	1:B:232:PRO:HG2	2.15	0.46
1:D:115:TYR:O	1:D:119:GLN:HG3	2.16	0.46
1:D:271:ILE:O	1:D:275:PHE:HB2	2.16	0.46
1:C:146:LEU:HB2	1:B:152:VAL:HG11	1.97	0.46
1:C:160:THR:HB	1:C:221:THR:O	2.15	0.46
1:B:256:VAL:HG12	1:B:257:PRO:HD3	1.97	0.46
1:D:149:THR:C	1:D:151:PRO:HD2	2.41	0.46
1:A:135:TYR:O	1:A:138:ILE:HG12	2.15	0.46
1:A:198:LEU:HB3	1:A:201:THR:OG1	2.15	0.46
1:B:36:GLU:OE2	1:B:40:GLU:HG3	2.15	0.46
1:A:102:CYS:CA	1:A:107:VAL:HG12	2.46	0.46
1:A:194:ASN:C	1:C:87:ARG:HD3	2.41	0.46
1:C:32:LYS:NZ	1:C:35:ARG:HG2	2.31	0.46
1:B:101:ASN:OD1	1:B:105:GLY:HA3	2.16	0.46
1:B:149:THR:C	1:B:151:PRO:HD2	2.40	0.46
1:B:194:ASN:C	1:D:87:ARG:HD3	2.41	0.46
1:A:219:MET:HG2	1:C:61:ASN:HB3	1.97	0.46
1:C:195:ASN:CB	1:C:196:PRO:HD3	2.45	0.46
1:B:32:LYS:C	1:B:35:ARG:HG2	2.40	0.46
1:B:33:MET:HE1	1:B:115:TYR:HE1	1.80	0.46
1:B:255:TRP:O	1:B:258:VAL:HG12	2.16	0.46
1:D:263:LEU:HD23	1:D:263:LEU:O	2.15	0.46
1:C:39:ALA:HA	1:C:42:MET:HG2	1.96	0.46
1:C:149:THR:OG1	1:C:151:PRO:CD	2.64	0.46
1:D:33:MET:HE1	1:D:115:TYR:CD1	2.50	0.46
1:B:48:MET:O	1:B:52:LEU:N	2.25	0.46
1:A:108:PRO:HG2	1:A:111:LYS:HD2	1.98	0.45
1:A:178:LEU:HD21	1:A:217:LEU:HD13	1.99	0.45
1:D:192:GLN:OE1	1:D:192:GLN:N	2.30	0.45
1:D:253:TRP:CZ3	1:D:256:VAL:HG11	2.50	0.45
1:A:102:CYS:SG	1:A:107:VAL:CG1	3.04	0.45
1:C:172:PHE:CZ	1:C:259:VAL:HG13	2.51	0.45
1:B:206:ILE:O	1:B:210:VAL:HG23	2.17	0.45
1:A:170:ARG:NH1	1:C:136:THR:HG23	2.26	0.45
1:C:82:VAL:CG1	1:C:203:ALA:HB2	2.46	0.45
1:D:271:ILE:HA	1:D:274:VAL:HG12	1.97	0.45
1:A:45:TYR:O	1:A:49:VAL:HG22	2.17	0.45
1:A:146:LEU:HD23	1:D:152:VAL:HA	1.98	0.45
1:B:173:LEU:HD22	1:D:133:LEU:HD23	1.98	0.45
1:D:72:LEU:HD11	1:D:76:PHE:CZ	2.51	0.45
1:D:130:ILE:HD11	1:D:134:PHE:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:184:LEU:HD12	1:D:271:ILE:HG13	1.98	0.45
1:B:33:MET:CE	1:B:114:VAL:HB	2.45	0.45
1:D:206:ILE:HD13	2:D:403:GOL:O1	2.15	0.45
1:D:216:SER:OG	1:D:217:LEU:HD12	2.16	0.45
1:D:228:SER:O	1:D:232:PRO:HG2	2.16	0.45
1:A:212:ILE:HD11	1:C:76:PHE:HB3	1.99	0.45
1:B:198:LEU:HD13	1:B:201:THR:CB	2.45	0.45
1:A:88:ILE:HD12	1:A:89:SER:CB	2.46	0.45
1:A:277:GLY:O	1:A:280:ILE:HG12	2.16	0.45
1:B:178:LEU:HD11	1:B:217:LEU:HB3	1.98	0.45
1:D:82:VAL:HG22	2:D:403:GOL:C1	2.46	0.45
1:C:276:ILE:HA	1:C:279:THR:HB	1.98	0.45
1:B:160:THR:HG21	1:B:224:ALA:HB2	1.99	0.45
1:B:164:ASP:N	1:B:164:ASP:OD1	2.47	0.45
1:B:274:VAL:HG13	1:B:275:PHE:HD1	1.80	0.45
1:A:108:PRO:HG2	1:A:111:LYS:CD	2.46	0.44
1:C:183:GLN:NE2	1:C:187:PHE:CZ	2.86	0.44
1:B:36:GLU:OE1	1:B:115:TYR:CE1	2.70	0.44
1:D:32:LYS:HD2	1:D:34:VAL:HG12	2.00	0.44
1:D:182:LEU:CD1	1:D:225:ILE:HG21	2.38	0.44
1:A:36:GLU:HB3	1:A:115:TYR:OH	2.17	0.44
1:A:182:LEU:HD22	1:A:225:ILE:HG21	1.99	0.44
1:B:32:LYS:O	1:B:35:ARG:HG2	2.18	0.44
1:B:212:ILE:O	1:B:215:VAL:HG12	2.17	0.44
1:D:60:LEU:HB3	1:D:137:ALA:HB1	2.00	0.44
1:A:33:MET:O	1:A:36:GLU:N	2.50	0.44
1:A:41:PHE:HZ	1:A:125:LEU:HD22	1.82	0.44
1:C:79:THR:HG22	1:C:80:MET:HE2	1.98	0.44
1:B:80:MET:O	1:B:84:VAL:HG12	2.17	0.44
1:A:39:ALA:O	1:A:42:MET:HG2	2.17	0.44
1:A:223:TYR:O	1:A:223:TYR:CD2	2.71	0.44
1:B:67:TYR:CZ	1:D:63:LYS:HD2	2.53	0.44
1:B:187:PHE:O	1:B:191:ASP:HB2	2.18	0.44
1:D:36:GLU:OE1	1:D:115:TYR:CE1	2.70	0.44
1:D:111:LYS:O	1:D:115:TYR:CD1	2.71	0.44
1:C:146:LEU:O	1:B:152:VAL:HG21	2.17	0.44
1:C:152:VAL:HB	1:B:146:LEU:HB3	1.98	0.44
1:C:230:ASP:O	1:C:234:ARG:HG3	2.18	0.44
1:C:268:GLY:O	1:C:271:ILE:HG22	2.17	0.44
1:D:160:THR:HB	1:D:221:THR:O	2.18	0.44
1:D:248:SER:HA	1:D:252:ASN:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:SER:OG	1:A:217:LEU:HD12	2.18	0.44
1:C:67:TYR:HE2	1:C:219:MET:HB2	1.82	0.44
1:B:41:PHE:HZ	1:B:125:LEU:HD22	1.83	0.44
1:B:162:LEU:HD11	1:B:166:MET:HG3	2.00	0.44
1:C:41:PHE:HZ	1:C:125:LEU:HD22	1.82	0.44
1:D:162:LEU:HD12	1:D:163:PRO:CD	2.47	0.44
1:A:194:ASN:OD1	1:C:87:ARG:CB	2.66	0.44
1:B:67:TYR:HE1	1:D:64:TYR:O	2.00	0.44
1:B:111:LYS:HD3	1:B:115:TYR:OH	2.17	0.44
1:B:253:TRP:O	1:B:253:TRP:CG	2.71	0.44
1:A:107:VAL:HG22	1:A:108:PRO:HD2	2.00	0.43
1:A:37:PHE:CZ	1:A:118:GLY:CA	3.01	0.43
1:A:168:LEU:HD13	1:A:252:ASN:O	2.19	0.43
1:A:80:MET:O	1:A:84:VAL:HG12	2.18	0.43
1:A:88:ILE:HD12	1:A:88:ILE:C	2.43	0.43
1:A:160:THR:HG21	1:A:224:ALA:HB2	2.00	0.43
1:C:149:THR:C	1:C:151:PRO:CD	2.91	0.43
1:C:237:THR:HG23	1:C:242:TRP:CE3	2.52	0.43
1:A:198:LEU:O	1:A:199:PRO:C	2.61	0.43
1:C:78:VAL:HG11	2:C:403:GOL:C1	2.48	0.43
1:C:163:PRO:HG3	1:C:219:MET:O	2.18	0.43
1:B:36:GLU:CD	1:B:115:TYR:CZ	2.97	0.43
1:D:168:LEU:HD11	1:D:252:ASN:HD22	1.84	0.43
1:A:162:LEU:HD11	1:A:166:MET:HG3	2.00	0.43
1:C:33:MET:HE1	1:C:115:TYR:HE1	1.79	0.43
1:C:45:TYR:CE1	1:C:126:ALA:HA	2.52	0.43
1:C:178:LEU:CD1	1:C:213:ILE:HG23	2.40	0.43
1:B:67:TYR:HE1	1:D:64:TYR:C	2.27	0.43
1:B:163:PRO:CD	1:B:220:ASN:O	2.66	0.43
1:A:116:VAL:HA	1:A:119:GLN:OE1	2.18	0.43
1:A:127:ALA:HA	1:A:130:ILE:HG22	2.01	0.43
1:A:166:MET:SD	1:A:220:ASN:O	2.76	0.43
1:A:195:ASN:O	1:A:196:PRO:C	2.59	0.43
1:C:72:LEU:HD11	1:C:76:PHE:CZ	2.53	0.43
1:B:67:TYR:CD1	1:D:63:LYS:HD2	2.53	0.43
1:B:163:PRO:HD3	1:B:220:ASN:O	2.18	0.43
1:D:150:GLY:N	1:D:151:PRO:HD3	2.34	0.43
1:A:160:THR:HG23	1:A:229:ARG:HH21	1.84	0.43
1:C:36:GLU:OE1	1:C:115:TYR:CE1	2.71	0.43
1:C:276:ILE:O	1:C:279:THR:HB	2.19	0.43
1:B:150:GLY:N	1:B:151:PRO:HD3	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:LEU:HD22	1:A:42:MET:CE	2.49	0.43
1:A:107:VAL:CG2	1:A:108:PRO:HD2	2.49	0.43
1:A:164:ASP:OD1	1:A:164:ASP:N	2.50	0.43
1:B:189:ILE:HG12	1:D:83:HIS:CE1	2.54	0.43
1:D:162:LEU:HD11	1:D:166:MET:HG3	2.00	0.43
1:A:111:LYS:HB3	1:A:115:TYR:CE2	2.53	0.43
1:C:164:ASP:OD1	1:C:164:ASP:N	2.51	0.43
1:B:256:VAL:HG13	1:B:257:PRO:HD3	2.01	0.43
1:D:37:PHE:CD2	1:D:114:VAL:O	2.72	0.43
1:D:52:LEU:HD23	1:D:130:ILE:HG21	2.01	0.43
1:A:152:VAL:HG11	1:D:146:LEU:CB	2.48	0.43
1:C:159:ALA:HA	1:C:229:ARG:HH12	1.84	0.43
1:C:94:ASN:HB3	1:C:97:VAL:HG12	2.00	0.42
1:B:161:TYR:CE1	1:B:249:ASN:HB2	2.54	0.42
1:B:232:PRO:HA	1:B:235:ILE:HG12	1.99	0.42
1:C:111:LYS:HB3	1:C:115:TYR:CE1	2.54	0.42
1:A:167:THR:HG21	1:A:169:TRP:CH2	2.54	0.42
1:A:267:LEU:O	1:A:267:LEU:HD23	2.19	0.42
1:C:80:MET:O	1:C:84:VAL:HG12	2.19	0.42
1:C:197:ALA:O	1:C:199:PRO:HD3	2.20	0.42
1:D:161:TYR:CE1	1:D:234:ARG:NH1	2.88	0.42
1:D:253:TRP:CG	1:D:253:TRP:O	2.71	0.42
1:A:93:MET:HE2	1:A:93:MET:N	2.35	0.42
1:A:115:TYR:O	1:A:119:GLN:OE1	2.38	0.42
1:C:182:LEU:HD11	2:C:403:GOL:H12	2.02	0.42
1:D:163:PRO:CD	1:D:220:ASN:O	2.67	0.42
1:D:237:THR:O	1:D:242:TRP:HB2	2.20	0.42
1:A:146:LEU:HD23	1:D:152:VAL:HG22	2.01	0.42
1:B:148:VAL:HG13	1:B:149:THR:N	2.35	0.42
1:B:166:MET:HE1	1:B:220:ASN:HB3	2.01	0.42
1:B:271:ILE:O	1:B:275:PHE:HB2	2.19	0.42
1:D:41:PHE:HZ	1:D:125:LEU:HD22	1.85	0.42
1:A:60:LEU:HB3	1:A:137:ALA:HB1	2.01	0.42
1:A:271:ILE:O	1:A:275:PHE:HB2	2.19	0.42
1:D:256:VAL:O	1:D:260:ALA:HB3	2.19	0.42
1:A:36:GLU:CD	1:A:115:TYR:OH	2.63	0.42
1:A:48:MET:HB3	1:A:126:ALA:CB	2.50	0.42
1:A:145:GLN:O	1:A:145:GLN:HG2	2.19	0.42
1:A:162:LEU:CG	1:A:166:MET:HG3	2.50	0.42
1:C:92:HIS:O	1:C:119:GLN:NE2	2.52	0.42
1:B:154:THR:O	1:B:157:ILE:HG12	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:LEU:HD11	1:B:166:MET:CG	2.50	0.42
1:B:184:LEU:HD12	1:B:271:ILE:HG13	2.02	0.42
1:D:45:TYR:CE1	1:D:126:ALA:HA	2.54	0.42
1:D:79:THR:HG22	1:D:80:MET:HE2	2.01	0.42
1:A:38:LEU:HD23	1:A:38:LEU:O	2.19	0.42
1:D:75:GLY:HA2	1:D:210:VAL:CG1	2.50	0.42
1:D:230:ASP:O	1:D:234:ARG:HG3	2.20	0.42
1:D:232:PRO:HA	1:D:235:ILE:HG12	2.01	0.42
1:A:189:ILE:HG23	1:A:198:LEU:HD12	2.02	0.42
1:C:127:ALA:HA	1:C:130:ILE:HG22	2.01	0.42
1:B:172:PHE:CZ	1:B:259:VAL:HG13	2.55	0.42
1:C:52:LEU:HD22	1:C:130:ILE:HG12	2.02	0.41
1:C:145:GLN:O	1:C:145:GLN:HG2	2.20	0.41
1:C:252:ASN:CG	1:C:252:ASN:O	2.59	0.41
1:B:195:ASN:O	1:B:196:PRO:C	2.62	0.41
1:B:279:THR:CG2	1:D:35:ARG:HG2	2.50	0.41
1:A:194:ASN:O	1:A:195:ASN:O	2.39	0.41
1:D:101:ASN:OD1	1:D:105:GLY:HA3	2.20	0.41
1:C:243:GLY:O	1:C:246:VAL:HG12	2.20	0.41
1:B:33:MET:HE3	1:B:114:VAL:HG11	2.03	0.41
1:D:185:CYS:CB	1:D:209:LEU:HD22	2.50	0.41
1:A:32:LYS:NZ	1:A:35:ARG:HB2	2.35	0.41
1:A:43:SER:HB3	1:A:93:MET:SD	2.61	0.41
1:D:260:ALA:HB3	1:D:261:PRO:HD3	2.02	0.41
1:A:37:PHE:CD2	1:A:114:VAL:O	2.74	0.41
1:A:105:GLY:O	1:A:106:ARG:HG2	2.20	0.41
1:B:37:PHE:CD2	1:B:114:VAL:O	2.74	0.41
1:B:198:LEU:HD21	1:D:83:HIS:HE1	1.85	0.41
1:D:162:LEU:HD13	1:D:253:TRP:CZ3	2.54	0.41
1:A:178:LEU:HD11	1:A:217:LEU:HB3	2.03	0.41
1:C:163:PRO:CG	1:C:219:MET:O	2.69	0.41
1:C:234:ARG:NH2	1:C:246:VAL:O	2.54	0.41
1:B:194:ASN:OD1	1:D:87:ARG:CB	2.69	0.41
1:D:34:VAL:HG22	1:D:34:VAL:O	2.19	0.41
1:A:149:THR:C	1:A:151:PRO:HD3	2.45	0.41
1:A:276:ILE:HD11	1:C:42:MET:SD	2.61	0.41
1:C:52:LEU:HD23	1:C:52:LEU:C	2.46	0.41
1:C:263:LEU:O	1:C:263:LEU:HD23	2.20	0.41
1:B:67:TYR:HE2	1:B:215:VAL:HG22	1.84	0.41
1:B:237:THR:O	1:B:242:TRP:HB2	2.21	0.41
1:D:164:ASP:OD1	1:D:164:ASP:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:THR:O	1:A:157:ILE:HG12	2.21	0.41
1:A:178:LEU:HD22	1:A:213:ILE:HG23	2.03	0.41
1:C:33:MET:HE2	1:C:114:VAL:CG2	2.51	0.41
1:B:223:TYR:O	1:B:223:TYR:CD2	2.74	0.41
1:C:253:TRP:CZ3	1:C:256:VAL:HG11	2.56	0.41
1:B:53:GLY:HA2	1:B:130:ILE:HD11	2.03	0.41
1:B:145:GLN:HG2	1:B:145:GLN:O	2.21	0.41
1:D:145:GLN:O	1:D:145:GLN:HG2	2.21	0.41
1:D:177:TRP:O	1:D:181:MET:HG3	2.20	0.41
1:A:53:GLY:HA2	1:A:130:ILE:HD11	2.02	0.41
1:C:44:THR:N	1:C:93:MET:SD	2.94	0.41
1:C:59:VAL:HG23	1:C:60:LEU:CG	2.47	0.41
1:B:33:MET:O	1:B:36:GLU:N	2.54	0.41
1:D:223:TYR:CD2	1:D:223:TYR:O	2.74	0.41
1:C:39:ALA:HB1	1:C:85:ALA:HB1	2.03	0.40
1:B:166:MET:SD	1:B:220:ASN:O	2.79	0.40
1:A:189:ILE:HG22	1:A:202:GLU:HG2	2.04	0.40
1:C:52:LEU:HD11	1:C:158:PHE:CD2	2.56	0.40
1:C:216:SER:OG	1:C:217:LEU:HD12	2.21	0.40
1:A:163:PRO:HG3	1:A:219:MET:O	2.21	0.40
1:D:48:MET:HB3	1:D:126:ALA:CB	2.51	0.40
1:D:185:CYS:HB2	1:D:209:LEU:HD22	2.03	0.40
1:D:207:GLY:O	1:D:210:VAL:HG12	2.21	0.40
1:A:44:THR:HG22	1:A:123:SER:OG	2.22	0.40
1:A:57:HIS:ND1	1:A:61:ASN:HB3	2.36	0.40
1:A:82:VAL:CG1	1:A:203:ALA:HB2	2.52	0.40
1:A:248:SER:O	1:A:251:GLU:O	2.39	0.40
1:C:57:HIS:O	1:C:63:LYS:HB3	2.22	0.40
1:C:104:LEU:HD11	1:C:273:LEU:HD22	2.02	0.40
1:B:93:MET:HE2	1:B:93:MET:N	2.35	0.40
1:C:166:MET:CE	1:C:220:ASN:HD22	2.34	0.40
1:B:225:ILE:O	1:B:261:PRO:HB3	2.21	0.40
1:D:38:LEU:O	1:D:38:LEU:HD23	2.22	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:VAL:O	1:C:194:ASN:HD21[26_555]	1.51	0.09
1:B:84:VAL:O	1:D:194:ASN:HD21[26_555]	1.51	0.09

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	247/376 (66%)	226 (92%)	21 (8%)	0	100	100
1	B	247/376 (66%)	228 (92%)	17 (7%)	2 (1%)	16	52
1	C	247/376 (66%)	230 (93%)	15 (6%)	2 (1%)	16	52
1	D	247/376 (66%)	228 (92%)	16 (6%)	3 (1%)	10	42
All	All	988/1504 (66%)	912 (92%)	69 (7%)	7 (1%)	18	54

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	252	ASN
1	B	253	TRP
1	D	252	ASN
1	D	253	TRP
1	C	195	ASN
1	B	195	ASN
1	D	195	ASN

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	195/304 (64%)	195 (100%)	0	100	100
1	B	195/304 (64%)	195 (100%)	0	100	100
1	C	195/304 (64%)	195 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	195/304 (64%)	195 (100%)	0	100	100
All	All	780/1216 (64%)	780 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	249	ASN
1	C	61	ASN
1	C	220	ASN
1	B	61	ASN
1	B	249	ASN
1	D	252	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	C	402	-	5,5,5	0.89	0	5,5,5	1.06	0
2	GOL	A	402	-	5,5,5	0.86	0	5,5,5	1.01	0
2	GOL	D	403	-	5,5,5	0.89	0	5,5,5	1.10	0
2	GOL	C	401	-	5,5,5	1.38	1 (20%)	5,5,5	0.90	0
2	GOL	B	403	-	5,5,5	0.98	0	5,5,5	1.07	0
2	GOL	D	402	-	5,5,5	0.89	0	5,5,5	1.09	0
2	GOL	D	404	-	5,5,5	0.88	0	5,5,5	1.04	0
2	GOL	B	402	-	5,5,5	0.89	0	5,5,5	1.02	0
2	GOL	C	403	-	5,5,5	1.02	0	5,5,5	0.86	0
2	GOL	A	401	-	5,5,5	1.19	0	5,5,5	1.04	0
2	GOL	D	401	-	5,5,5	1.33	1 (20%)	5,5,5	1.14	1 (20%)
2	GOL	B	401	-	5,5,5	1.42	2 (40%)	5,5,5	0.84	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	C	402	-	-	2/4/4/4	-
2	GOL	A	402	-	-	0/4/4/4	-
2	GOL	D	403	-	-	0/4/4/4	-
2	GOL	C	401	-	-	2/4/4/4	-
2	GOL	B	403	-	-	0/4/4/4	-
2	GOL	D	402	-	-	0/4/4/4	-
2	GOL	D	404	-	-	2/4/4/4	-
2	GOL	B	402	-	-	0/4/4/4	-
2	GOL	C	403	-	-	2/4/4/4	-
2	GOL	A	401	-	-	2/4/4/4	-
2	GOL	D	401	-	-	0/4/4/4	-
2	GOL	B	401	-	-	0/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	GOL	C1-C2	2.43	1.61	1.51
2	D	401	GOL	C1-C2	2.24	1.60	1.51
2	B	401	GOL	C1-C2	2.24	1.60	1.51
2	B	401	GOL	C3-C2	2.19	1.60	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	GOL	C3-C2-C1	-2.09	104.13	111.80

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	402	GOL	O1-C1-C2-C3
2	C	403	GOL	O1-C1-C2-O2
2	C	403	GOL	O1-C1-C2-C3
2	A	401	GOL	O2-C2-C3-O3
2	D	404	GOL	O1-C1-C2-O2
2	A	401	GOL	C1-C2-C3-O3
2	D	404	GOL	O1-C1-C2-C3
2	C	402	GOL	O1-C1-C2-O2
2	C	401	GOL	O1-C1-C2-C3
2	C	401	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	403	GOL	7	0
2	C	403	GOL	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	249/376 (66%)	1.05	26 (10%)	11 14	86, 97, 105, 127	0
1	B	249/376 (66%)	0.74	13 (5%)	33 27	68, 80, 87, 92	0
1	C	249/376 (66%)	0.98	30 (12%)	9 12	83, 98, 105, 122	0
1	D	249/376 (66%)	0.72	18 (7%)	21 20	71, 80, 89, 106	0
All	All	996/1504 (66%)	0.88	87 (8%)	16 17	68, 90, 103, 127	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	36	GLU	8.1
1	C	195	ASN	6.3
1	A	252	ASN	5.3
1	A	35	ARG	4.8
1	D	195	ASN	4.5
1	B	252	ASN	4.2
1	C	268	GLY	4.1
1	D	252	ASN	4.0
1	D	180	GLY	3.8
1	C	252	ASN	3.8
1	A	86	GLY	3.7
1	A	40	GLU	3.6
1	C	180	GLY	3.6
1	A	115	TYR	3.5
1	A	67	TYR	3.5
1	B	194	ASN	3.4
1	C	63	LYS	3.4
1	D	268	GLY	3.4
1	A	194	ASN	3.3
1	B	170	ARG	3.2
1	D	196	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	63	LYS	3.2
1	A	268	GLY	3.1
1	C	271	ILE	3.1
1	C	226	ASN	3.1
1	A	207	GLY	3.1
1	C	196	PRO	3.0
1	B	229	ARG	3.0
1	A	57	HIS	2.9
1	A	243	GLY	2.9
1	C	194	ASN	2.9
1	C	35	ARG	2.8
1	C	248	SER	2.8
1	C	280	ILE	2.8
1	D	57	HIS	2.8
1	A	51	GLY	2.7
1	D	102	CYS	2.7
1	A	227	PRO	2.6
1	B	180	GLY	2.6
1	C	102	CYS	2.6
1	D	280	ILE	2.6
1	B	171	GLY	2.6
1	B	243	GLY	2.6
1	C	278	SER	2.5
1	D	248	SER	2.5
1	D	168	LEU	2.5
1	D	194	ASN	2.5
1	A	199	PRO	2.5
1	A	170	ARG	2.5
1	C	253	TRP	2.4
1	B	57	HIS	2.4
1	D	260	ALA	2.4
1	B	67	TYR	2.4
1	B	268	GLY	2.4
1	C	168	LEU	2.4
1	C	260	ALA	2.4
1	A	171	GLY	2.3
1	C	142	SER	2.3
1	A	142	SER	2.3
1	A	180	GLY	2.3
1	C	86	GLY	2.3
1	C	259	VAL	2.3
1	A	200	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	277	GLY	2.3
1	C	32	LYS	2.3
1	A	151	PRO	2.3
1	C	221	THR	2.2
1	C	170	ARG	2.2
1	C	243	GLY	2.2
1	D	243	GLY	2.2
1	B	211	VAL	2.2
1	C	157	ILE	2.2
1	C	237	THR	2.2
1	D	221	THR	2.2
1	C	40	GLU	2.2
1	C	277	GLY	2.2
1	C	100	ALA	2.1
1	A	259	VAL	2.1
1	A	260	ALA	2.1
1	A	258	VAL	2.1
1	B	280	ILE	2.1
1	A	119	GLN	2.1
1	A	195	ASN	2.0
1	C	171	GLY	2.0
1	D	253	TRP	2.0
1	B	260	ALA	2.0
1	D	203	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	D	404	6/6	0.43	0.24	74,88,90,91	14
2	GOL	D	403	6/6	0.67	0.53	73,88,89,89	14
2	GOL	D	402	6/6	0.76	0.18	79,95,98,99	14
2	GOL	C	402	6/6	0.80	0.15	82,99,110,111	14
2	GOL	B	401	6/6	0.83	0.25	95,116,133,139	0
2	GOL	C	401	6/6	0.84	0.19	112,134,157,159	0
2	GOL	A	401	6/6	0.85	0.18	110,132,144,145	0
2	GOL	B	403	6/6	-	-	71,85,87,88	14
2	GOL	C	403	6/6	0.85	0.52	89,91,109,109	14
2	GOL	A	402	6/6	0.86	0.11	78,94,101,102	14
2	GOL	B	402	6/6	0.88	0.11	69,82,87,89	14
2	GOL	D	401	6/6	0.90	0.20	97,122,141,147	0

6.5 Other polymers [i](#)

There are no such residues in this entry.